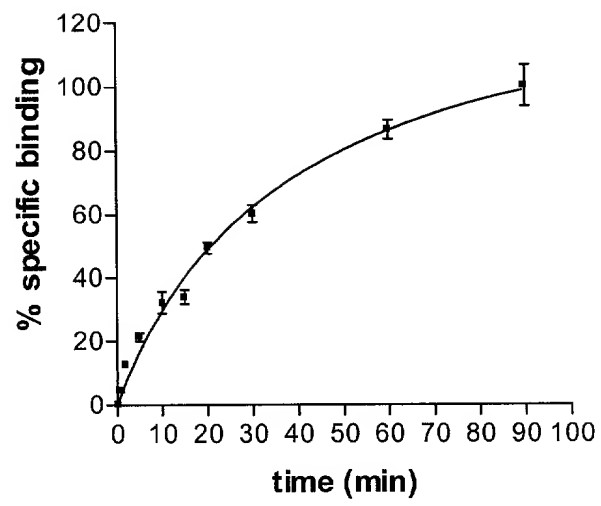


Figure 1: Association [3H]-4MG



**Figure 2: Dissociation of
[3H]-4MG binding**

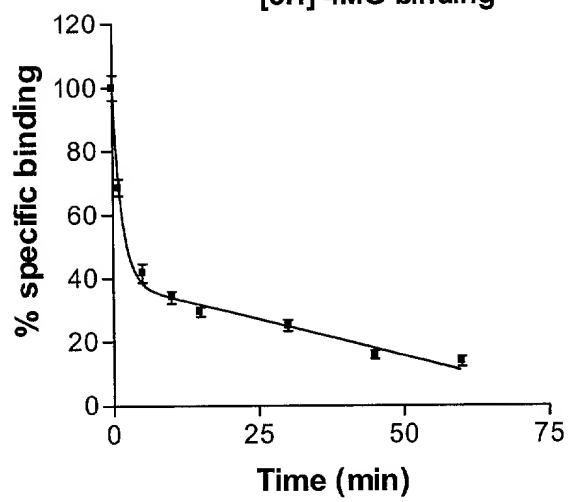


Figure 3: DRUG INHIBITION OF
[3H]-4MG BINDING

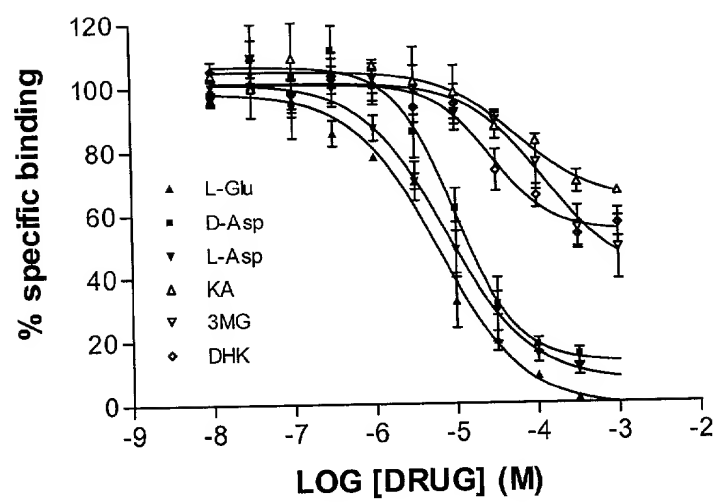


Figure 4: DRUG INHIBITION of [3H]-4MG
BINDING

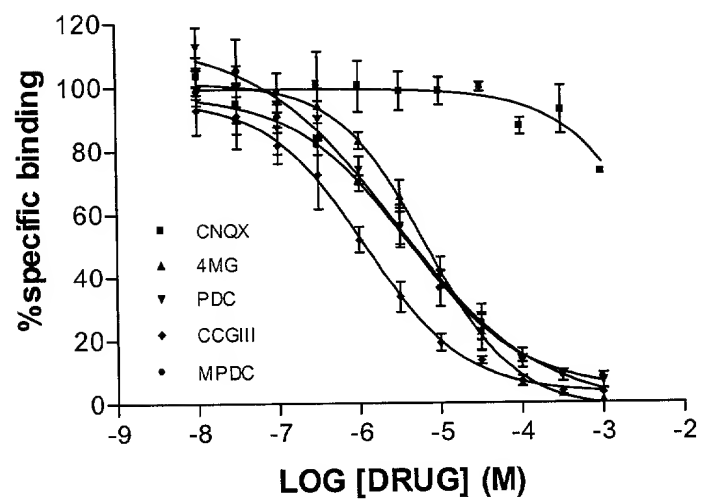
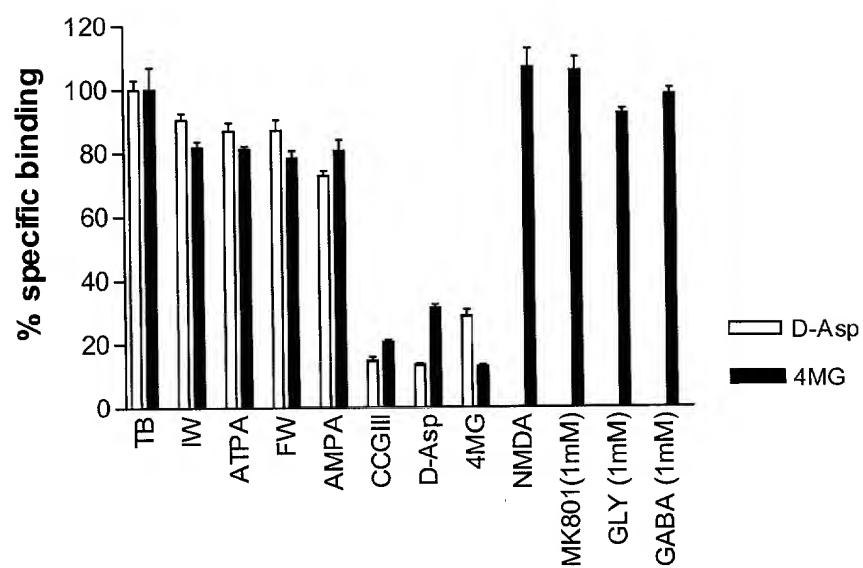


Figure 5 : COMPARISON of MISCELLANEOUS DRUG
INHIBITION of [3H]-D-ASPARTATE and [3H]-4MG



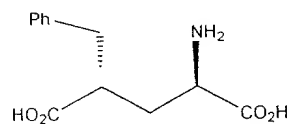
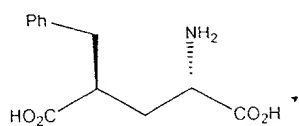
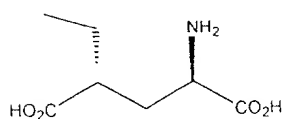
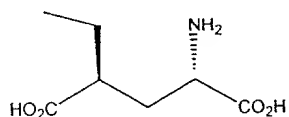
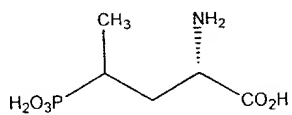
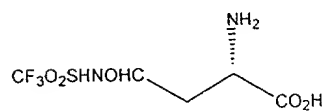
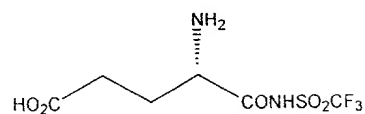
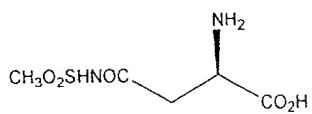
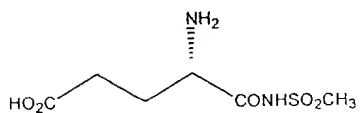
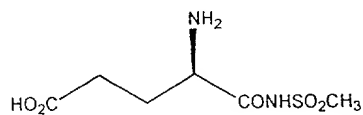
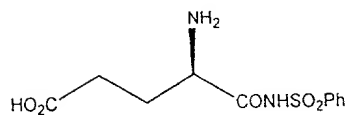
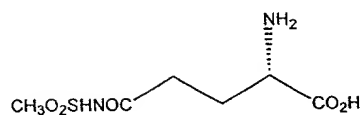
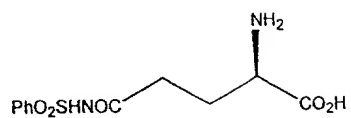
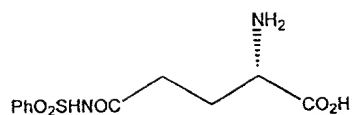
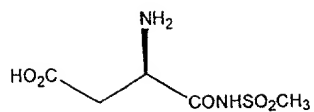
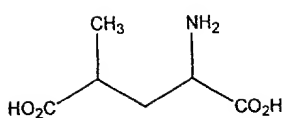


Figure 6A

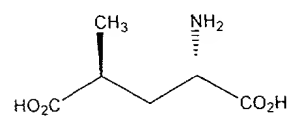
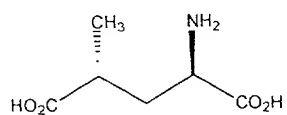
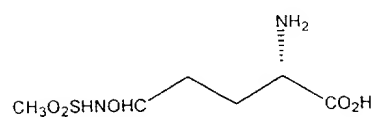
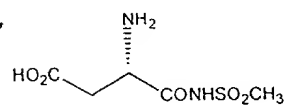
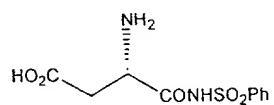
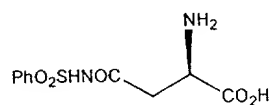
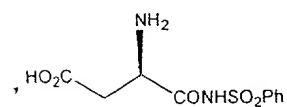
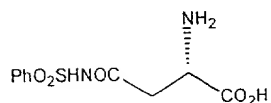
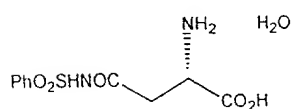
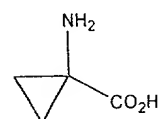
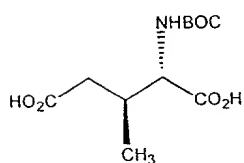
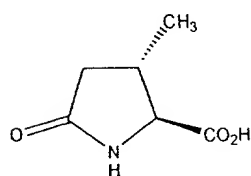
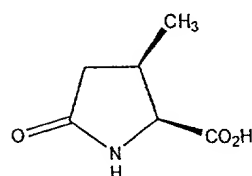
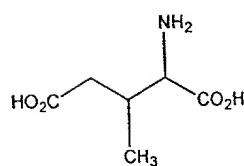
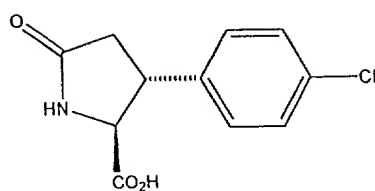
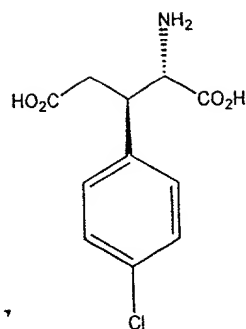
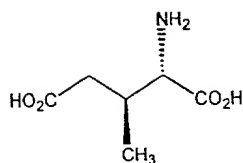
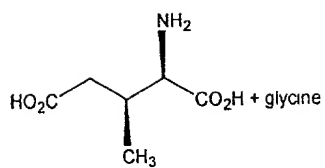


Figure 6B

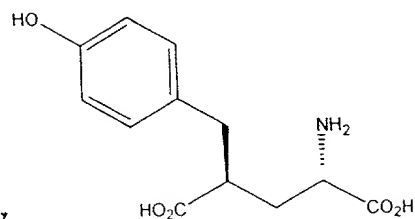
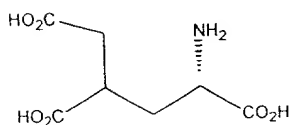
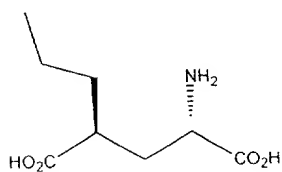
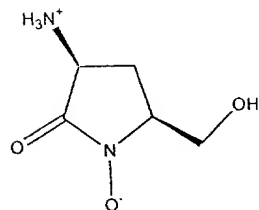
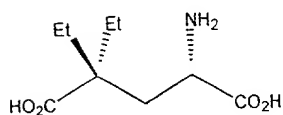
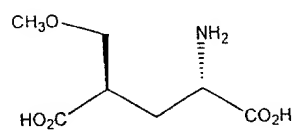
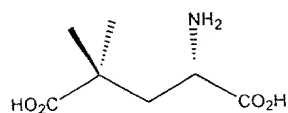
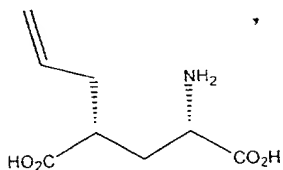
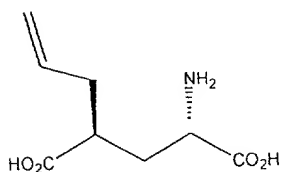
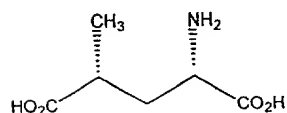
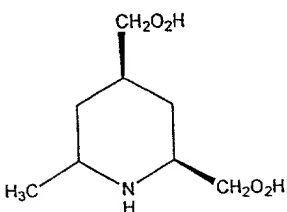
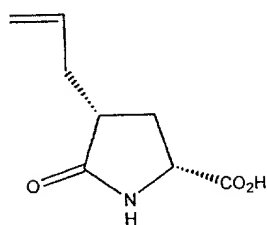
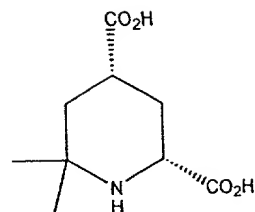
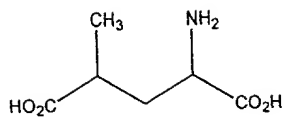
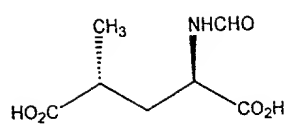


Figure 6C

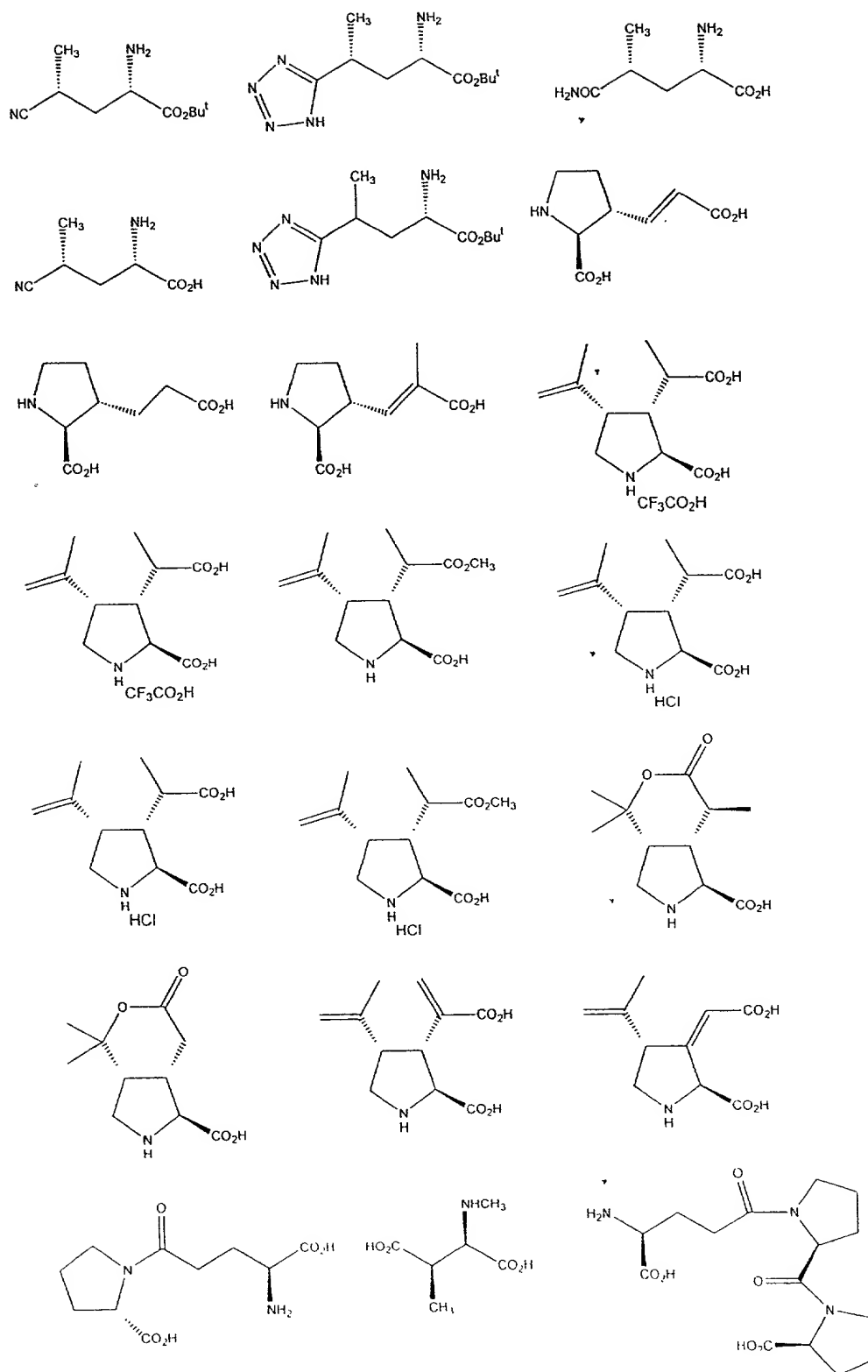


Figure 6E

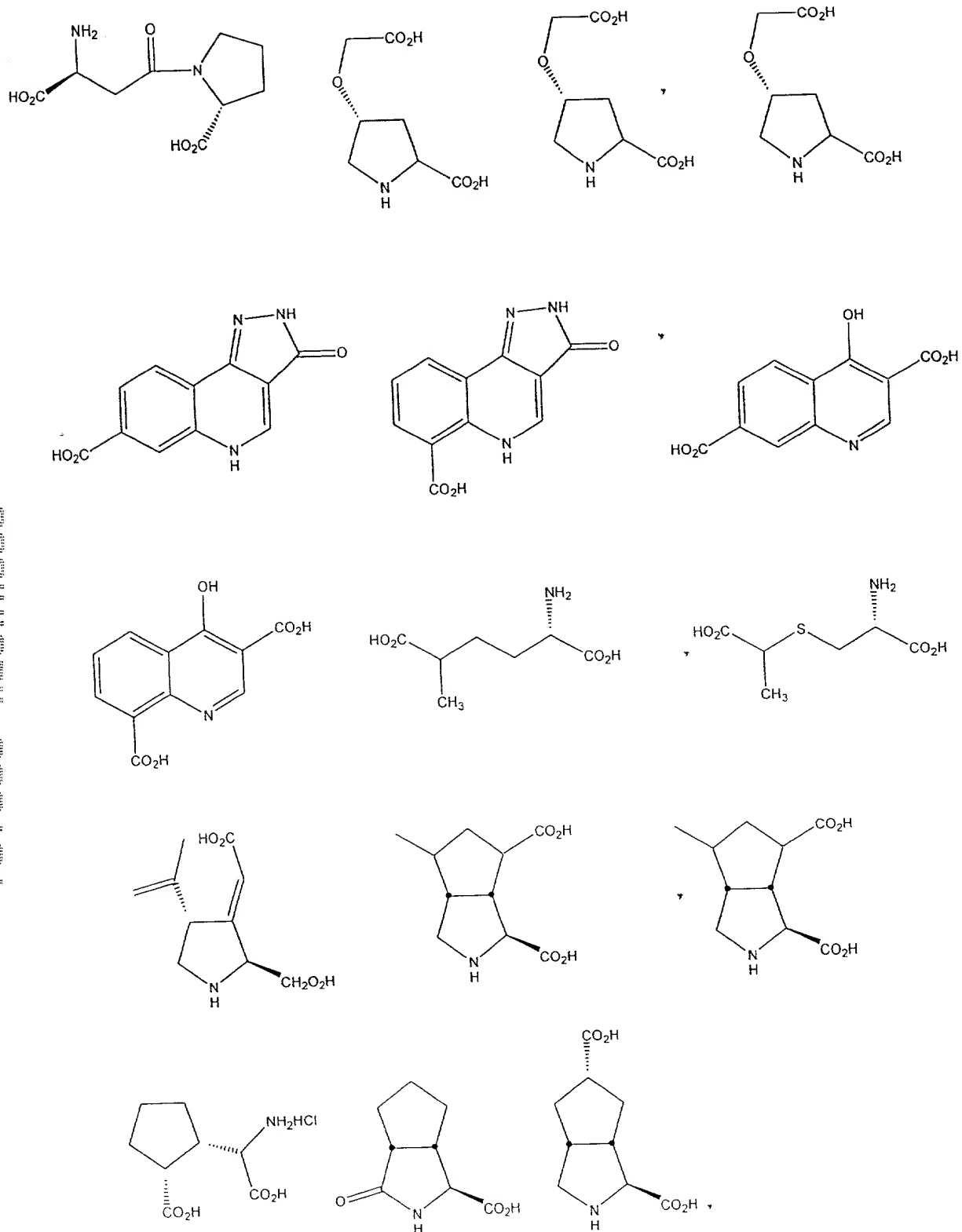


Figure 6F

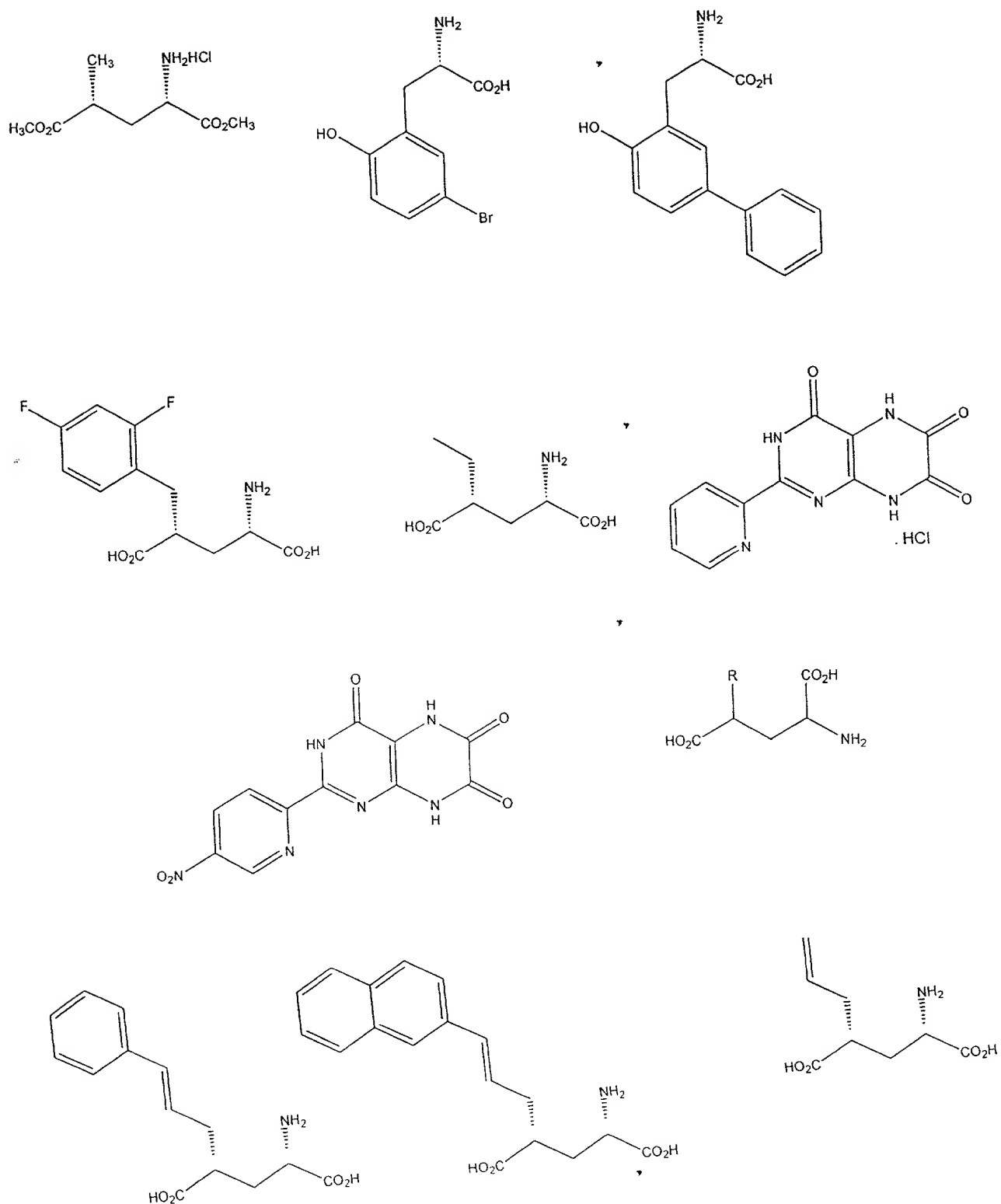


Figure 6G

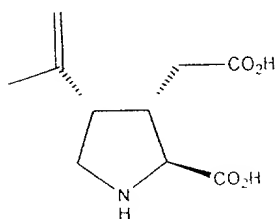
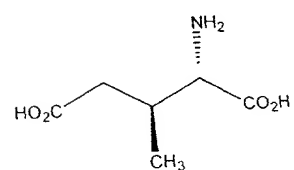
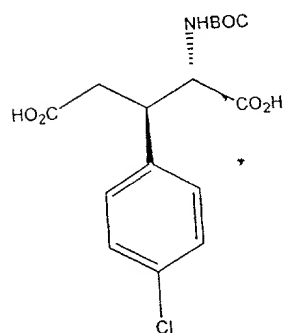
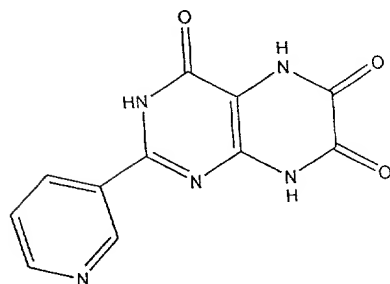
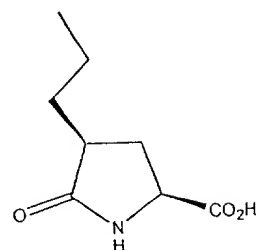
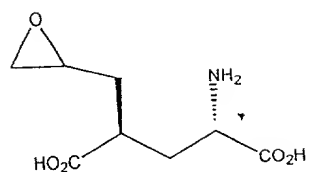
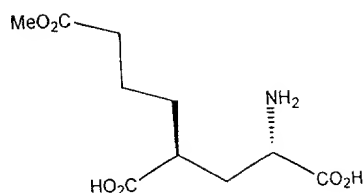
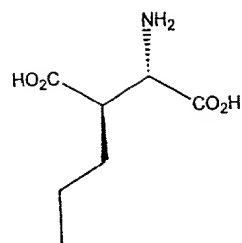
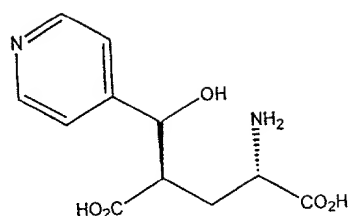
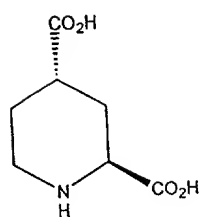
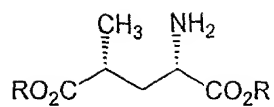
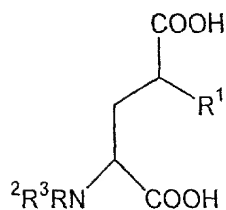
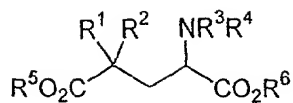


Figure 6H



R = Me, Et, ^tBu

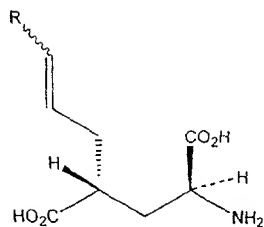
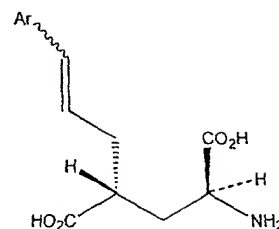


R¹ = CH₃, and halogen

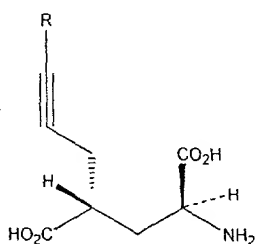
R², R³ are independently

H, C1-C6-alkyl, C3-C4-alkenyl, C3-C5-cycloalkyl, C1-C6-alkyl-CO-,
C1-C6-alkyl-OCO-, C1-C6-alkyl-NHCO-, HCO-, or C3-C6-alkynyl

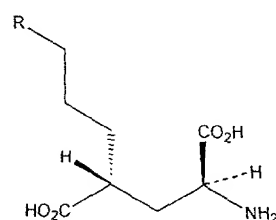
R², R³ taken together can be -CH₂(CH₂)_pCH₂-



R = H, Me, Et, Cl



R = H, Me, Et, nPr



R = H, Et, nPr

Figure 6I